



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 13, 2026 – 10:54 PM UTC

PDB ID : 7CBZ / pdb_00007cbz
Title : Crystal structure of T2R-TTL-A31 complex
Authors : Yang, J.H.; Yan, W.
Deposited on : 2020-06-15
Resolution : 2.61 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtrriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

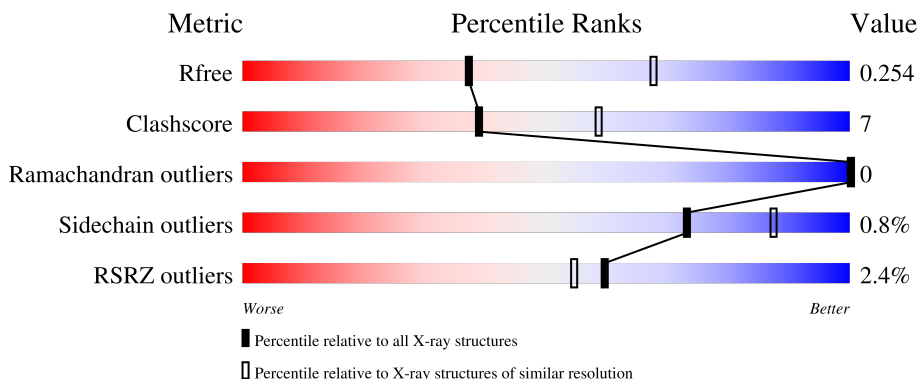
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.61 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



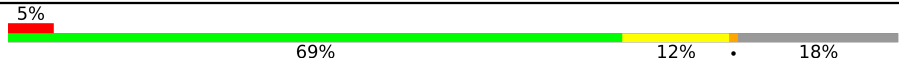
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4951 (2.64-2.60)
Clashscore	190562	5303 (2.64-2.60)
Ramachandran outliers	187476	5217 (2.64-2.60)
Sidechain outliers	187428	5217 (2.64-2.60)
RSRZ outliers	180081	4950 (2.64-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	451	
1	C	451	
2	B	445	
2	D	445	
3	E	189	

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Mol	Chain	Length	Quality of chain
4	F	383	 <p>5% 69% 12% 18%</p>

2 Entry composition [i](#)

There are 11 unique types of molecules in this entry. The entry contains 34298 atoms, of which 16837 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Tubulin alpha-1B chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	438	6769	2170	3341	582	654	22	0	1	0
1	C	440	6836	2190	3380	584	659	23	0	5	0

- Molecule 2 is a protein called Tubulin beta chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
2	B	424	6565	2101	3222	570	646	26	0	1	0
2	D	425	6553	2095	3215	569	648	26	0	1	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	57	THR	ALA	conflict	UNP P02554
B	172	MET	VAL	conflict	UNP P02554
B	298	SER	ALA	conflict	UNP P02554
B	318	ILE	VAL	conflict	UNP P02554
D	57	THR	ALA	conflict	UNP P02554
D	172	MET	VAL	conflict	UNP P02554
D	298	SER	ALA	conflict	UNP P02554
D	318	ILE	VAL	conflict	UNP P02554

- Molecule 3 is a protein called Stathmin-4.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
3	E	123	2085	638	1049	188	205	5	0	3	0

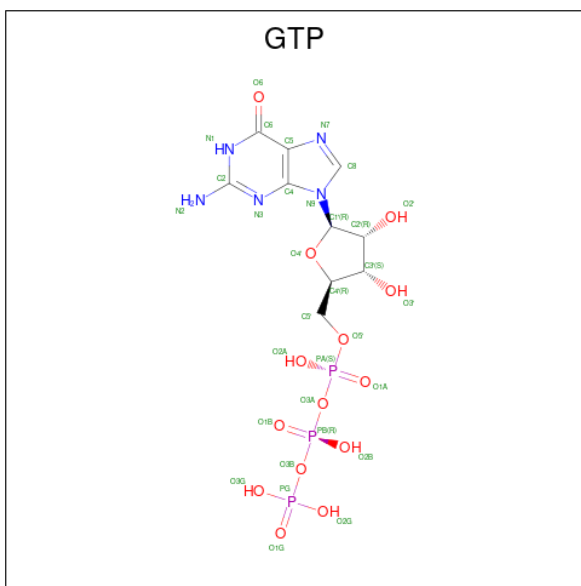
- Molecule 4 is a protein called Tubulin tyrosine ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
4	F	313	5032	1626	2507	425	462	12	0	0	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	379	HIS	-	expression tag	UNP E1BQ43
F	380	HIS	-	expression tag	UNP E1BQ43
F	381	HIS	-	expression tag	UNP E1BQ43
F	382	HIS	-	expression tag	UNP E1BQ43
F	383	HIS	-	expression tag	UNP E1BQ43

- Molecule 5 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: C₁₀H₁₆N₅O₁₄P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	H	N	O			P
5	A	1	41	10	9	5	14	3	41	0
5	C	1	41	10	9	5	14	3	41	0

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total 1 Mg 1	1	0
6	B	1	Total 1 Mg 1	1	0

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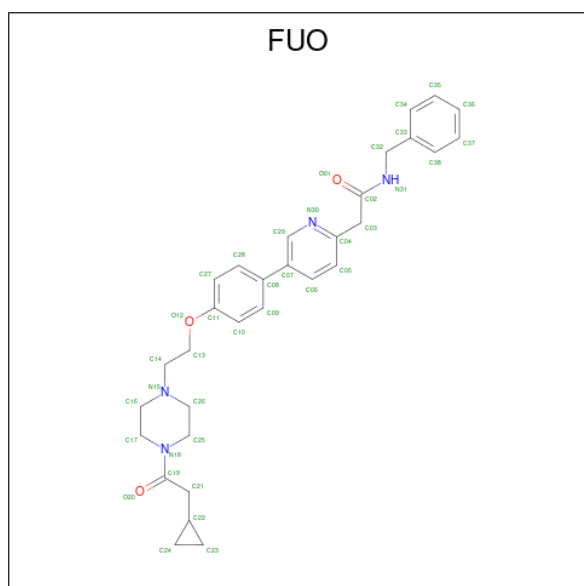
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	C	1	Total Mg 1 1	1	0
6	D	1	Total Mg 1 1	1	0
6	F	1	Total Mg 1 1	1	0

- Molecule 7 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	2	Total Ca 2 2	2	0
7	B	2	Total Ca 2 2	2	0
7	C	1	Total Ca 1 1	1	0
7	D	1	Total Ca 1 1	1	0

- Molecule 8 is 2-[5-[4-[2-[4-(2-cyclopropylethanoyl)piperazin-1-yl]ethoxy]phenyl]pyridin-2-yl]-N-(phenylmethyl)ethanamide (CCD ID: FUI) (formula: C₃₁H₃₆N₄O₃) (labeled as "Ligand of Interest" by depositor).



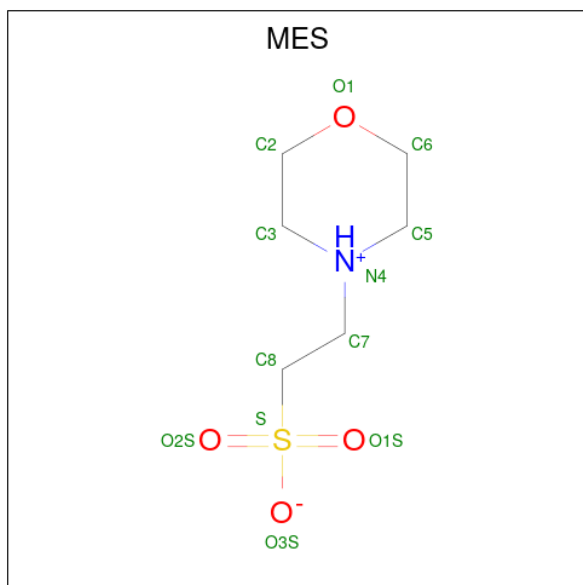
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	B	1	Total C H N O 74 31 36 4 3	0	0

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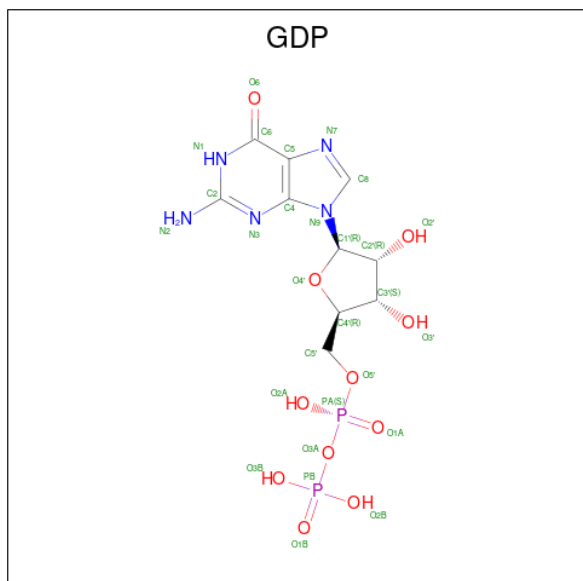
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	H	N	O		
8	D	1	74	31	36	4	3	0	0

- Molecule 9 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	H	N	O	S		
9	B	1	25	6	13	1	4	1	25	0

- Molecule 10 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula: $C_{10}H_{15}N_5O_{11}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	H	N	O			P
10	B	1	38	10	10	5	11	2	38	0
10	D	1	38	10	10	5	11	2	38	0

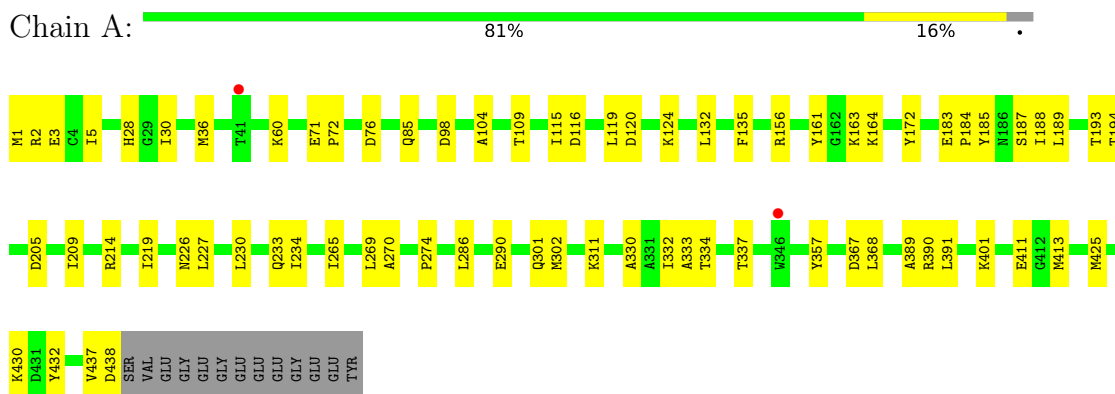
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	21	Total	O	0	0
			21	21		
11	B	35	Total	O	0	0
			35	35		
11	C	27	Total	O	0	0
			27	27		
11	D	14	Total	O	0	0
			14	14		
11	E	9	Total	O	0	0
			9	9		
11	F	10	Total	O	0	0
			10	10		

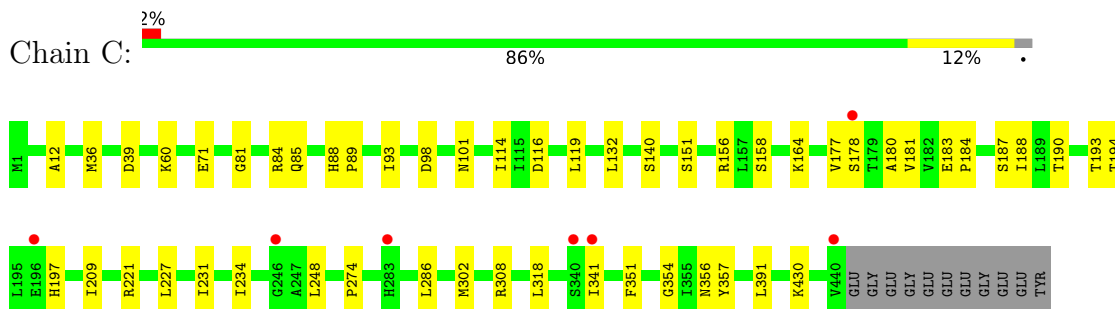
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

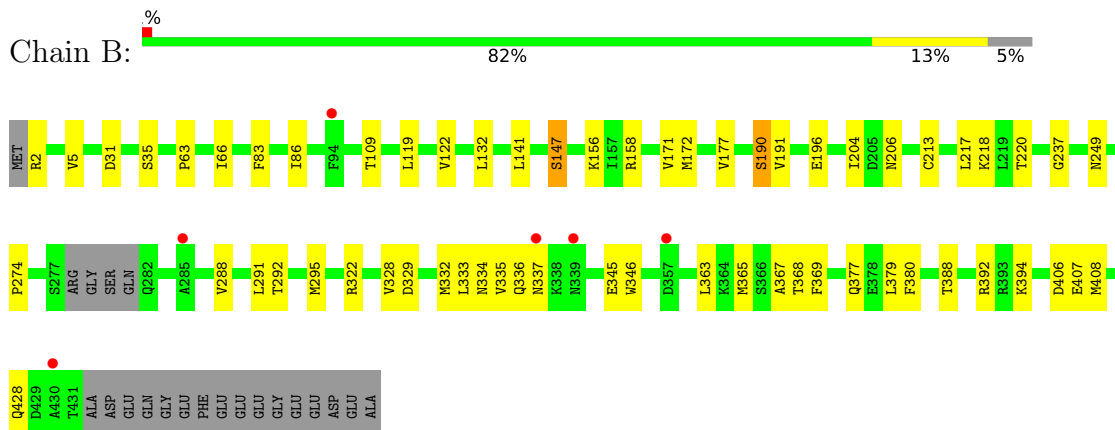
- Molecule 1: Tubulin alpha-1B chain



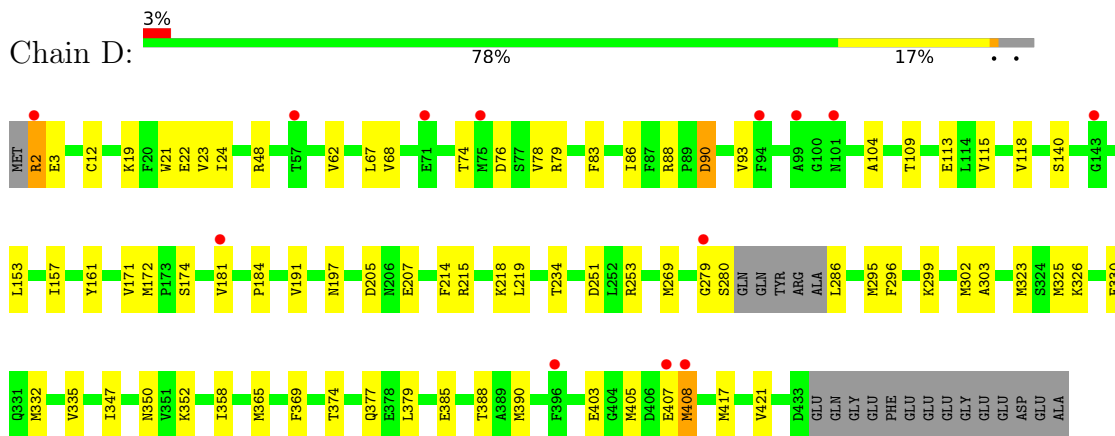
- Molecule 1: Tubulin alpha-1B chain



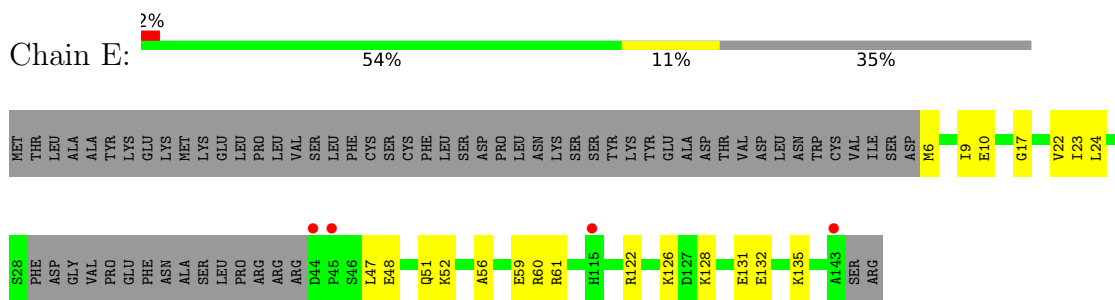
- Molecule 2: Tubulin beta chain



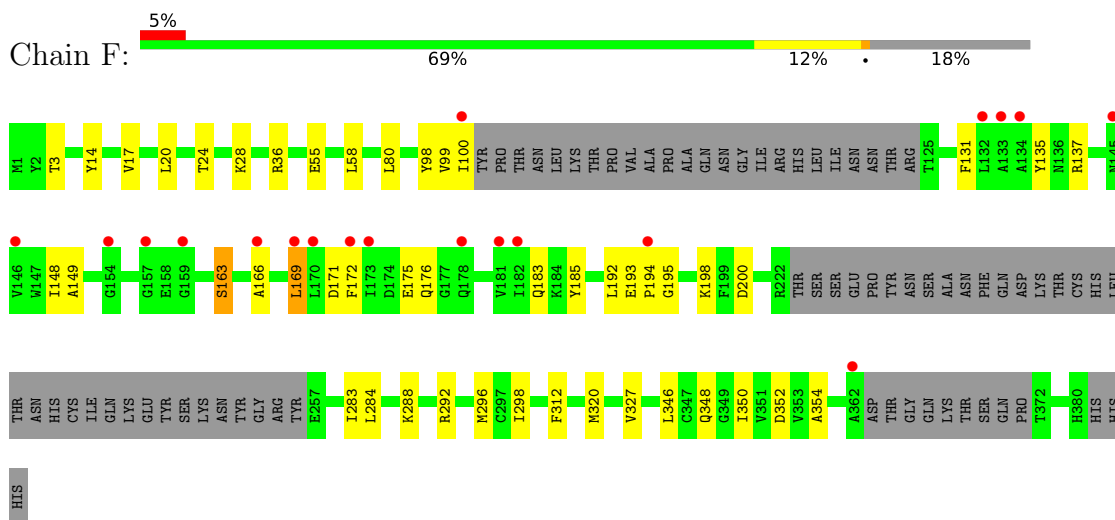
- Molecule 2: Tubulin beta chain



- Molecule 3: Stathmin-4



- Molecule 4: Tubulin tyrosine ligase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	104.81Å 156.63Å 181.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.76 – 2.61 19.76 – 2.61	Depositor EDS
% Data completeness (in resolution range)	92.9 (19.76-2.61) 92.9 (19.76-2.61)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.95 (at 2.59Å)	Xtrriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.206 , 0.255 0.207 , 0.254	Depositor DCC
R_{free} test set	4262 reflections (4.64%)	wwPDB-VP
Wilson B-factor (Å ²)	58.4	Xtrriage
Anisotropy	0.211	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 47.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	34298	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FUU, MES, GDP, GTP, CA, MG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.21	0/3509	0.41	0/4764
1	C	0.33	1/3549 (0.0%)	0.42	1/4819 (0.0%)
2	B	0.52	2/3420 (0.1%)	0.55	3/4633 (0.1%)
2	D	0.42	1/3414 (0.0%)	0.51	4/4624 (0.1%)
3	E	0.61	1/1048 (0.1%)	0.68	2/1391 (0.1%)
4	F	0.27	0/2578	0.50	2/3482 (0.1%)
All	All	0.39	5/17518 (0.0%)	0.49	12/23713 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	10	GLU	CA-C	-10.77	1.43	1.53
2	B	190	SER	C-O	-5.84	1.17	1.24
2	B	191	VAL	C-O	-5.53	1.17	1.24
1	C	356	ASN	C-O	-5.32	1.17	1.24
2	D	3	GLU	C-O	-5.15	1.18	1.24

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	147	SER	CA-C-N	7.36	127.97	119.94
2	B	147	SER	C-N-CA	7.36	127.97	119.94
1	C	354	GLY	O-C-N	-6.54	117.62	123.37
2	D	197	ASN	N-CA-C	6.43	121.17	112.88
2	D	251	ASP	O-C-N	-5.88	114.31	121.83

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3428	3341	3341	48	3
1	C	3456	3380	3380	40	3
2	B	3343	3222	3221	43	0
2	D	3338	3215	3215	56	0
3	E	1036	1049	1047	13	0
4	F	2525	2507	2506	40	0
5	A	32	9	12	0	0
5	C	32	9	12	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
6	F	1	0	0	0	0
7	A	2	0	0	0	0
7	B	2	0	0	0	0
7	C	1	0	0	0	0
7	D	1	0	0	0	0
8	B	38	36	0	0	0
8	D	38	36	0	2	0
9	B	12	13	13	0	0
10	B	28	10	12	0	0
10	D	28	10	12	0	0
11	A	21	0	0	2	0
11	B	35	0	0	4	0
11	C	27	0	0	0	0
11	D	14	0	0	1	0
11	E	9	0	0	1	0
11	F	10	0	0	0	0
All	All	17461	16837	16771	232	3

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

The worst 5 of 232 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:248:LEU:CD1	1:C:357:TYR:OH	1.99	1.11
1:C:248:LEU:HD12	1:C:357:TYR:OH	1.51	1.09
2:D:172:MET:HE2	2:D:379:LEU:HD21	1.45	0.98
1:A:290:GLU:OE2	11:A:601:HOH:O	1.89	0.89
4:F:163:SER:HB2	4:F:169:LEU:CD2	2.03	0.88

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:290:GLU:OE1	1:C:308:ARG:NH2[3_555]	1.59	0.61
1:A:290:GLU:OE1	1:C:308:ARG:HH21[3_555]	1.09	0.51
1:A:290:GLU:CD	1:C:308:ARG:NH2[3_555]	2.05	0.15

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	437/451 (97%)	425 (97%)	12 (3%)	0	100	100
1	C	443/451 (98%)	434 (98%)	9 (2%)	0	100	100
2	B	421/445 (95%)	405 (96%)	16 (4%)	0	100	100
2	D	422/445 (95%)	408 (97%)	14 (3%)	0	100	100
3	E	122/189 (65%)	120 (98%)	2 (2%)	0	100	100
4	F	305/383 (80%)	291 (95%)	14 (5%)	0	100	100
All	All	2150/2364 (91%)	2083 (97%)	67 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	370/379 (98%)	370 (100%)	0	100	100
1	C	376/379 (99%)	375 (100%)	1 (0%)	86	94
2	B	368/383 (96%)	365 (99%)	3 (1%)	73	87
2	D	368/383 (96%)	362 (98%)	6 (2%)	55	78
3	E	113/171 (66%)	110 (97%)	3 (3%)	39	65
4	F	273/341 (80%)	271 (99%)	2 (1%)	76	89
All	All	1868/2036 (92%)	1853 (99%)	15 (1%)	73	87

5 of 15 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	253	ARG
4	F	163	SER
2	D	407	GLU
4	F	169	LEU
3	E	128	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 28 such sidechains are listed below:

Mol	Chain	Res	Type
2	D	197	ASN
4	F	333	ASN
2	D	336	GLN
4	F	136	ASN
2	D	331	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 11 are monoatomic - leaving 7 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	GTP	A	501	6	33,34,34	0.92	1 (3%)	50,54,54	1.73	10 (20%)
8	FUO	D	501	-	42,42,42	3.40	18 (42%)	54,56,56	1.86	10 (18%)
9	MES	B	503	-	12,12,12	2.32	1 (8%)	15,16,16	1.55	3 (20%)
5	GTP	C	501	6	33,34,34	0.96	1 (3%)	50,54,54	1.64	9 (18%)
8	FUO	B	501	-	42,42,42	3.42	19 (45%)	54,56,56	2.90	17 (31%)
10	GDP	B	506	6	29,30,30	1.16	3 (10%)	45,47,47	1.83	7 (15%)
10	GDP	D	504	6	29,30,30	1.13	3 (10%)	45,47,47	1.97	10 (22%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GTP	A	501	6	-	8/22/38/38	0/3/3/3
8	FUO	D	501	-	-	6/27/39/39	0/5/5/5
9	MES	B	503	-	-	1/6/14/14	0/1/1/1
5	GTP	C	501	6	-	4/22/38/38	0/3/3/3
8	FUO	B	501	-	-	5/27/39/39	1/5/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
10	GDP	B	506	6	-	2/16/32/32	0/3/3/3
10	GDP	D	504	6	-	4/16/32/32	0/3/3/3

The worst 5 of 46 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	D	501	FUO	C37-C38	9.05	1.54	1.38
8	D	501	FUO	C34-C33	7.88	1.54	1.38
9	B	503	MES	C8-S	-7.76	1.66	1.77
8	B	501	FUO	C14-N15	-7.55	1.30	1.47
8	D	501	FUO	C36-C35	7.12	1.53	1.38

The worst 5 of 66 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	B	501	FUO	C03-C02-N31	13.80	135.21	116.06
8	B	501	FUO	O01-C02-N31	-7.56	108.19	123.03
10	D	504	GDP	C5-C4-N3	-6.58	117.92	128.39
10	B	506	GDP	C5-C4-N3	-6.29	118.38	128.39
8	D	501	FUO	C03-C02-N31	6.00	124.39	116.06

There are no chirality outliers.

5 of 30 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	501	GTP	C5'-O5'-PA-O3A
5	A	501	GTP	C5'-O5'-PA-O1A
5	A	501	GTP	C5'-O5'-PA-O2A
5	C	501	GTP	C5'-O5'-PA-O3A
5	C	501	GTP	C5'-O5'-PA-O2A

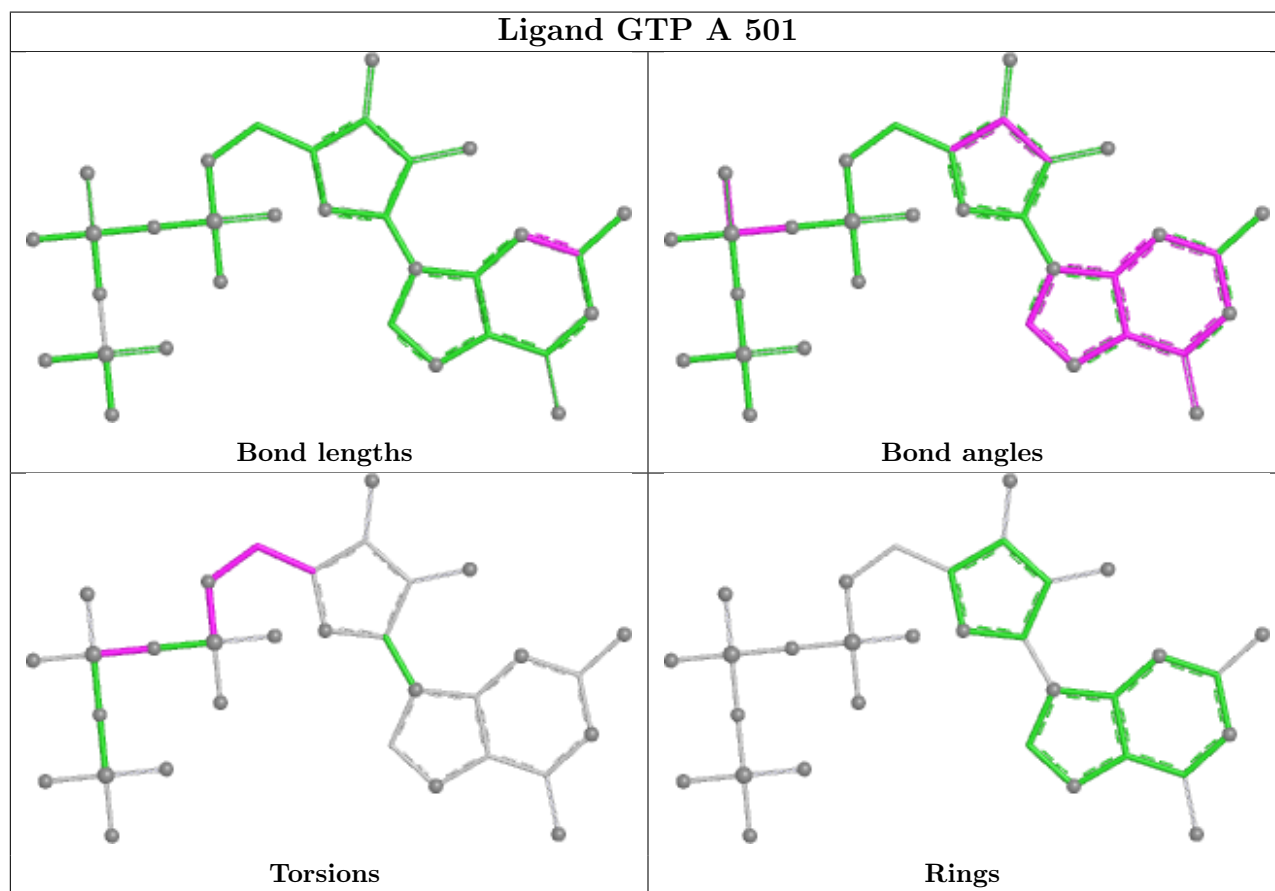
All (1) ring outliers are listed below:

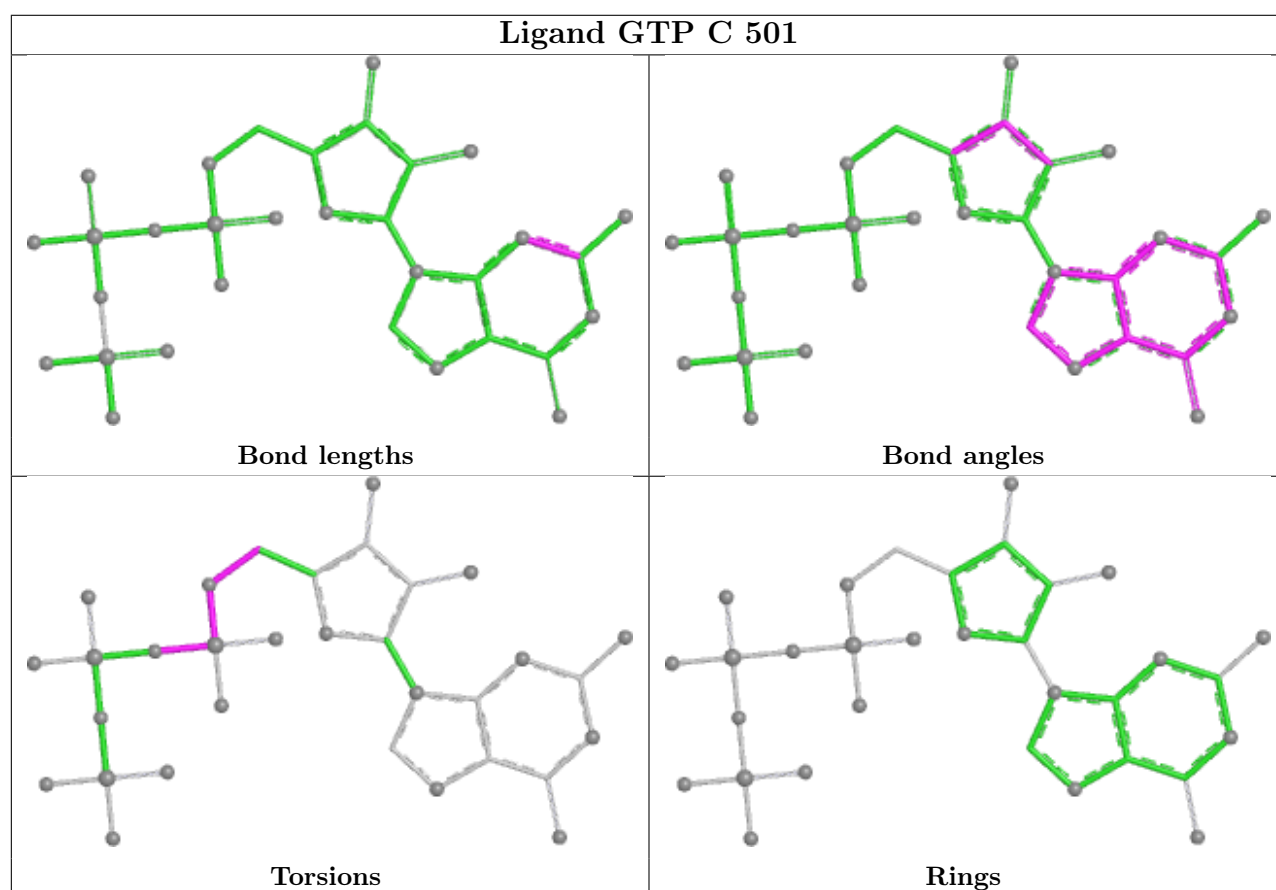
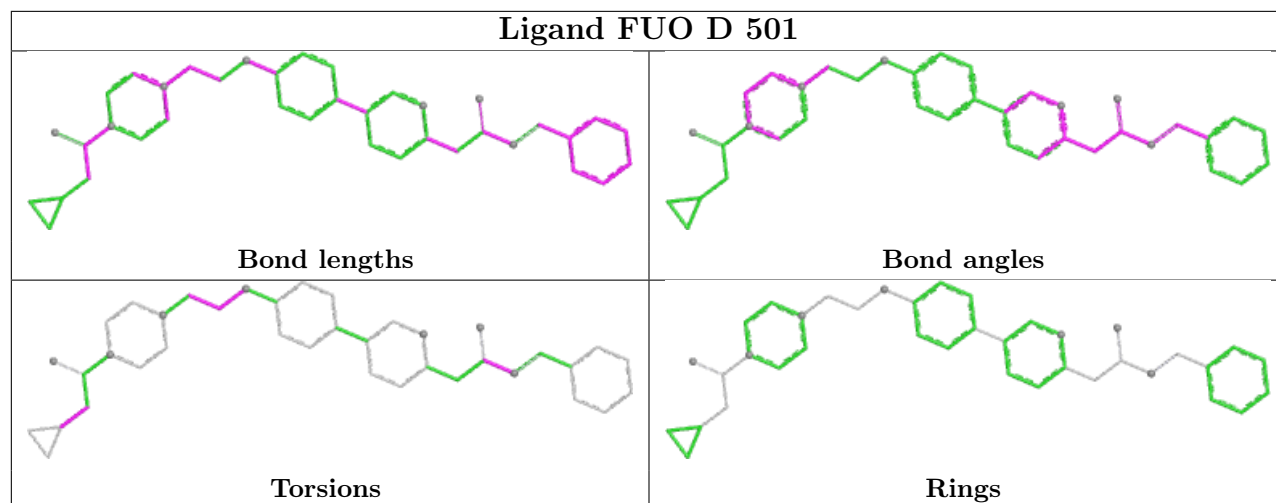
Mol	Chain	Res	Type	Atoms
8	B	501	FUO	C16-C17-C25-C26-N15-N18

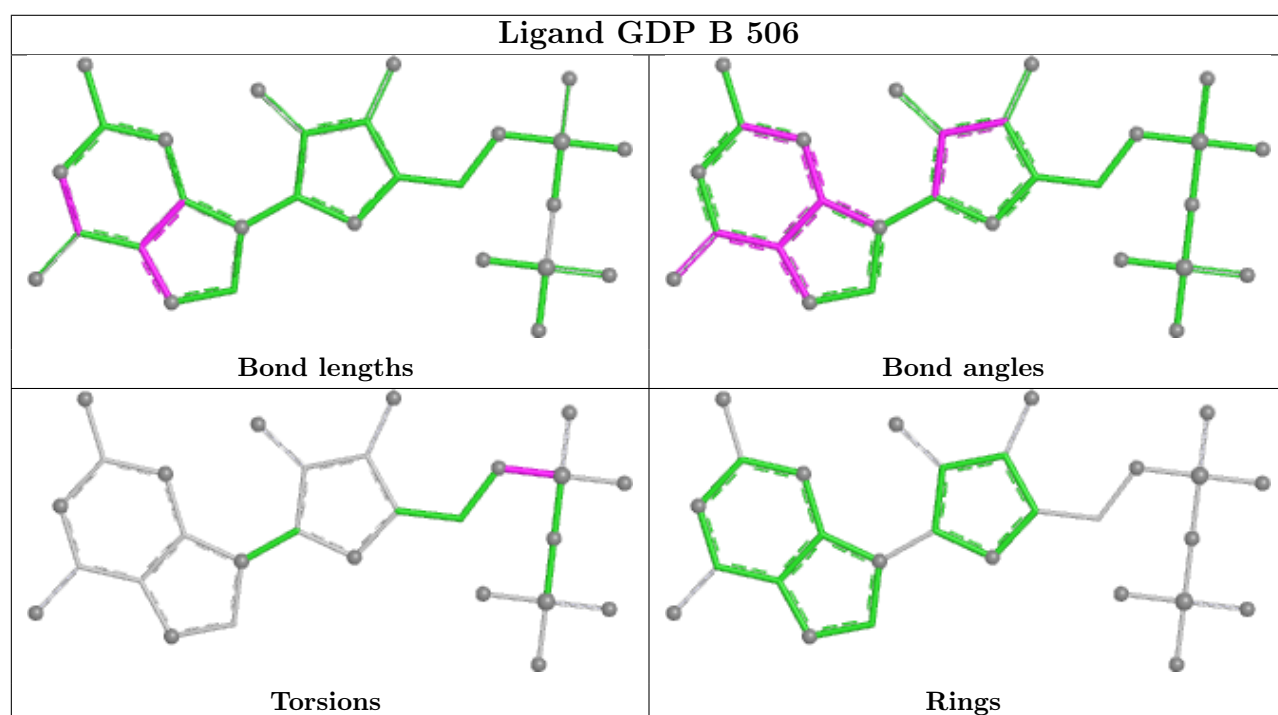
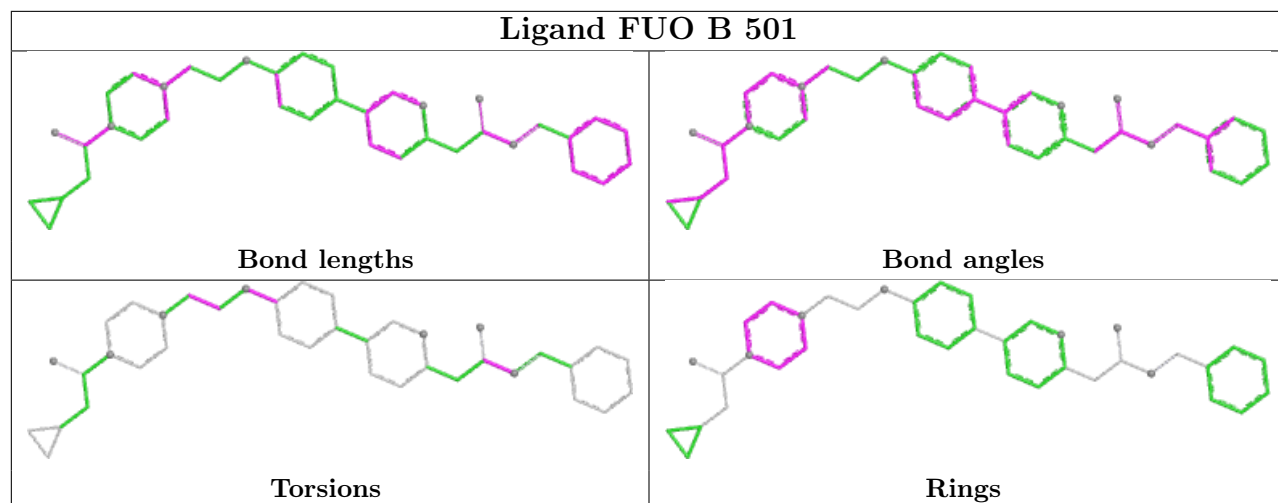
1 monomer is involved in 2 short contacts:

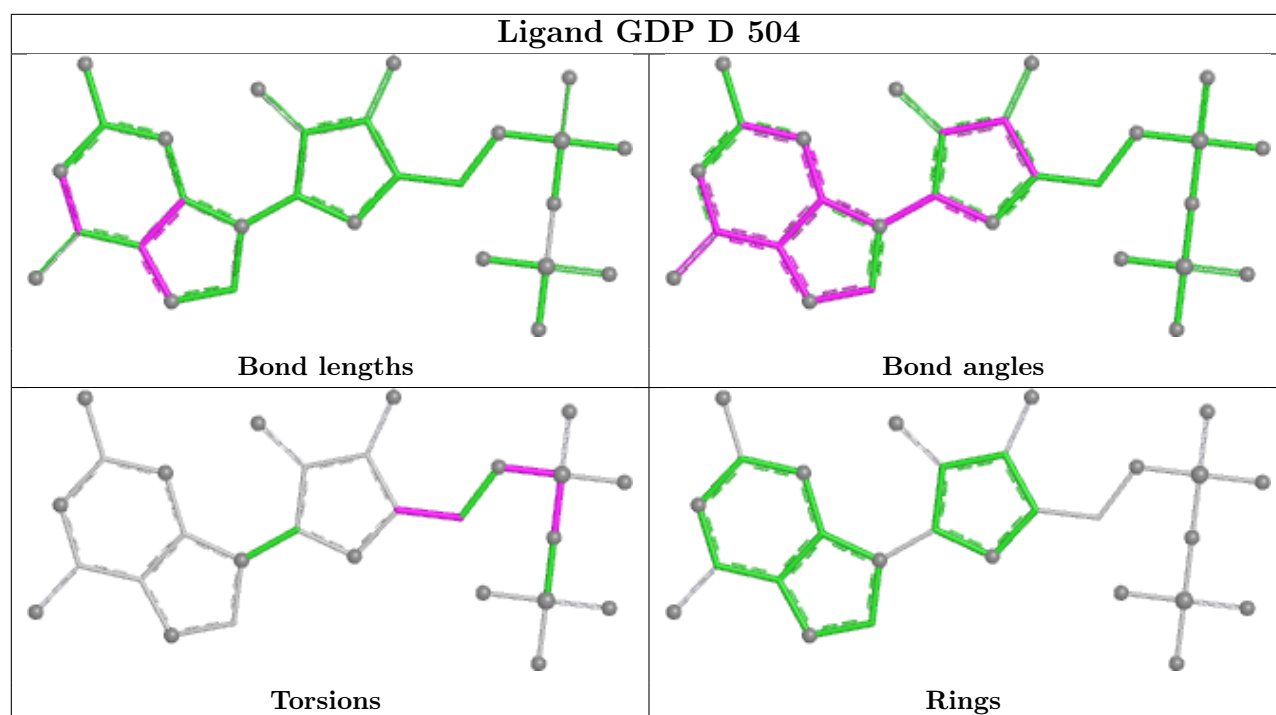
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	D	501	FUO	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	438/451 (97%)	-0.05	2 (0%) 87 85	37, 65, 96, 120	1 (0%)
1	C	440/451 (97%)	-0.27	7 (1%) 70 67	26, 55, 88, 125	5 (1%)
2	B	424/445 (95%)	-0.09	6 (1%) 73 70	36, 63, 99, 131	1 (0%)
2	D	425/445 (95%)	0.26	13 (3%) 51 46	40, 80, 113, 131	1 (0%)
3	E	123/189 (65%)	0.42	4 (3%) 49 44	29, 83, 118, 137	3 (2%)
4	F	313/383 (81%)	0.47	19 (6%) 27 22	51, 89, 158, 182	0
All	All	2163/2364 (91%)	0.06	51 (2%) 59 54	26, 69, 117, 182	11 (0%)

The worst 5 of 51 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	F	166	ALA	4.4
3	E	143	ALA	4.0
4	F	157	GLY	3.5
4	F	100	ILE	3.4
2	B	430	ALA	3.4

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

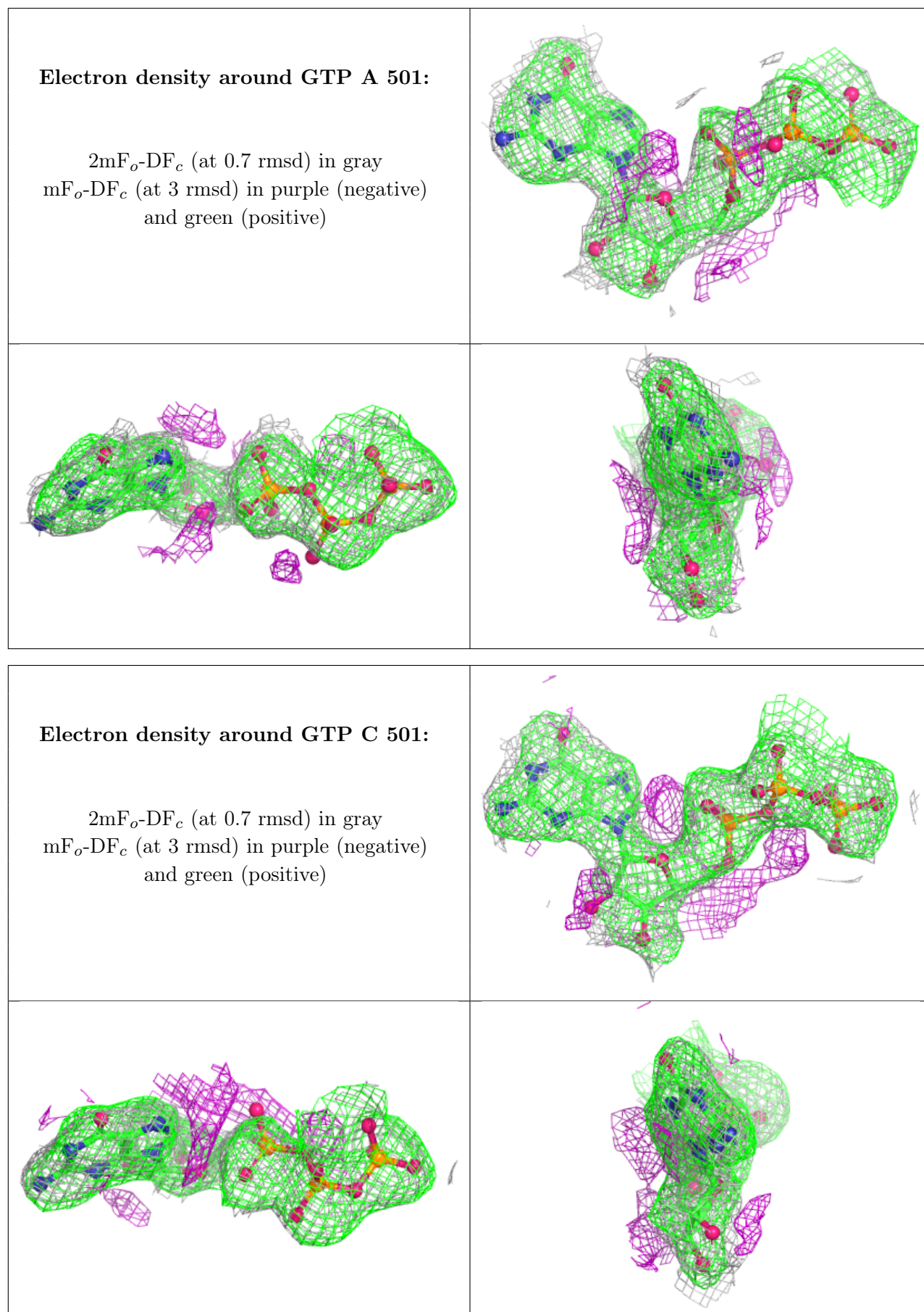
There are no oligosaccharides in this entry.

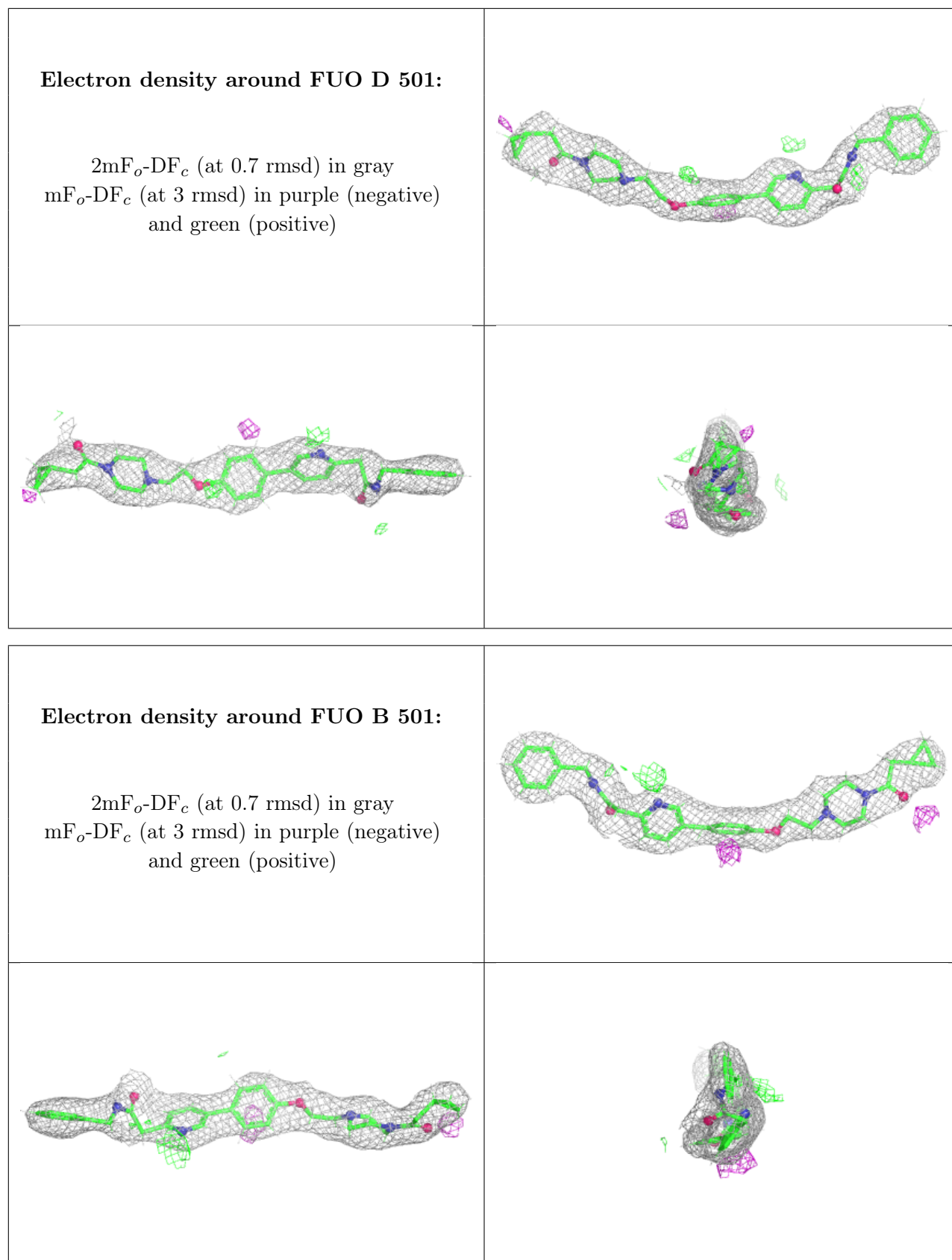
6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	GTP	A	501	32/32	-	-	44,46,56,57	41
5	GTP	C	501	32/32	-	-	38,42,51,53	41
6	MG	A	502	1/1	-	-	48,48,48,48	1
6	MG	B	502	1/1	-	-	50,50,50,50	1
6	MG	C	502	1/1	-	-	43,43,43,43	1
6	MG	D	502	1/1	-	-	150,150,150,150	1
6	MG	F	401	1/1	-	-	85,85,85,85	1
7	CA	A	503	1/1	-	-	82,82,82,82	1
7	CA	A	504	1/1	-	-	59,59,59,59	1
7	CA	B	504	1/1	-	-	83,83,83,83	1
7	CA	B	505	1/1	-	-	77,77,77,77	1
7	CA	C	503	1/1	-	-	62,62,62,62	1
7	CA	D	503	1/1	-	-	75,75,75,75	1
8	FUO	D	501	38/38	0.90	0.12	47,74,104,129	0
8	FUO	B	501	38/38	0.92	0.09	41,63,87,90	0
9	MES	B	503	12/12	-	-	46,49,57,57	25
10	GDP	B	506	28/28	-	-	40,45,55,56	38
10	GDP	D	504	28/28	-	-	83,108,152,173	38

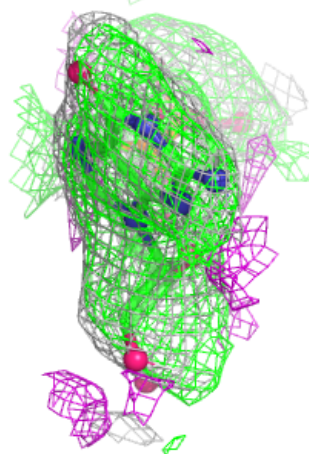
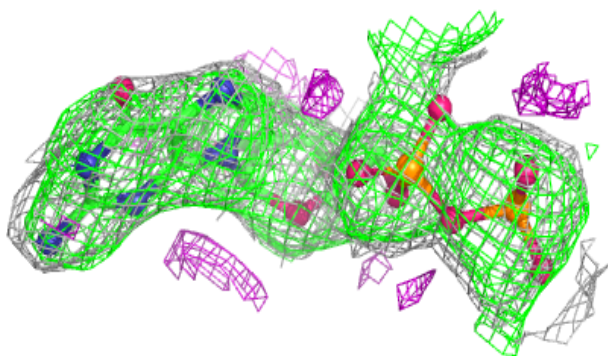
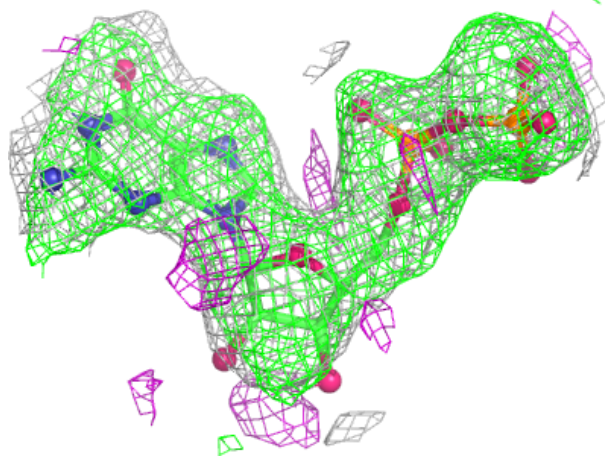
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

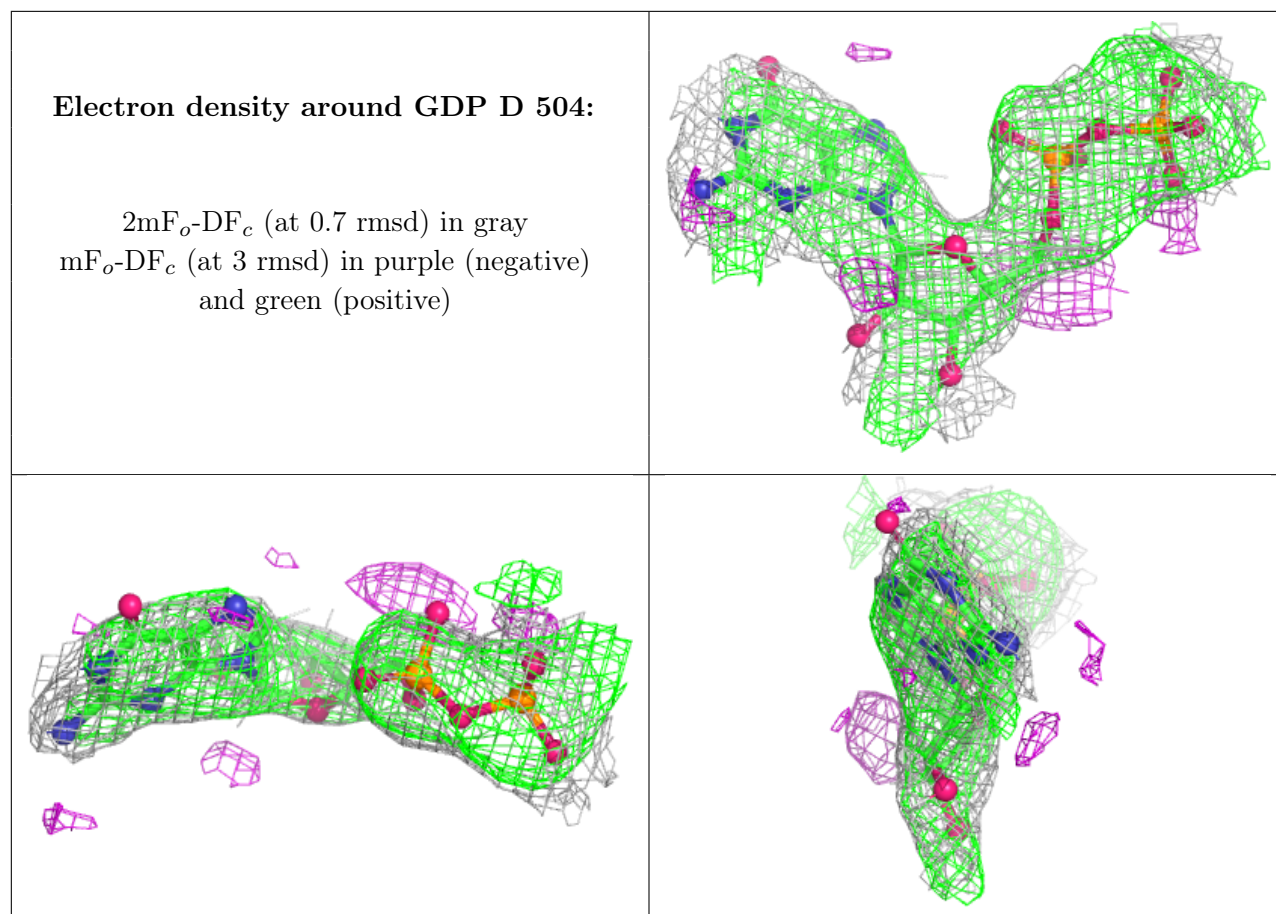




Electron density around GDP B 506:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers [i](#)

There are no such residues in this entry.